

Chemical and Environmental Measurement Information

# Recra LabNet Philadelphia Analytical Report \*\*REVISION\*\*

Client: TNU-HANFORD B99-085

**W.O.** #: 10985-001-001-9999-00

RFW#: 9908L851

Date Received: 08-24-99

SDG/SAF #: H0500/B99-085

#### **SEMIVOLATILE**

This narrative was corrected to add the TIC search for Tributylphosphate.

One (1) water sample was collected on 08-19-99.



**FDMC** 

The sample and its associated QC samples were extracted on 08-26-99 and analyzed according to criteria set forth in Recra OPs based on SW 846 Method 8270B for TCL Semivolatile target compounds on 09-13-99.

The following is a summary of the QC results accompanying the sample results and a description of any problems encountered during their analyses:

- 1. The cooler temperature upon receipt has been recorded on the chain-of-custody.
- 2. The required holding times for extraction and analysis were met.
- 3. Non-target compounds were detected in the sample.
- 4. All surrogate recoveries were within EPA QC limits.
- 5. One (1) of twenty-two (22) matrix spike recoveries was outside EPA QC limits.
- 6. All blank spike recoveries were within EPA QC limits.
- 7. The method blank contained the common laboratory contaminant Bis(2-Ethylhexyl)phthalate at a level greater than 5x the CRQL; however, no other compounds were detected in the samples. Re-extraction was not required. A copy of the Sample Discrepancy Report (SDR) has been enclosed.
- 8. A spectral search was conducted for the compound Butylated Hydroxytoluene and Tributylphosphate; these compounds were not identified in the samples.

J. Michael Taylor

Vice President

Date

pel/gorupsdata/bnastnu08851.doc

Philadelphia Analytical Laboratory

The results presented in this report relate only to the analytical testing and conditions of the samples at receipt and during storage. All pages of this report are integral parts of the analytical data. Therefore, this report should only be reproduced in its entirety of 10 pages, including page # 3A.

#### **GLOSSARY OF BNA DATA**

#### DATA QUALIFIERS

U	=	Compound was analyzed for but not detected. The associated numerical value is the estimated
		sample quantitation limit which is included and corrected for dilution and percent moisture.

- J = Indicates an estimated value. This flag is used under the following circumstances: 1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; or 2) when the mass spectral data indicate the presence of a compound that meets the identification criteria but the result is less than the specified detection limit but greater than zero. For example, if the limit of detection is 10 ug/L and a concentration of 3 ug/L is calculated, it is reported as 3J.
- B = This flag is used when the analyte is found in the associated blank as well as in the sample. It indicates possible/probable blank contamination. This flag is also used for a TIC as well as for a positively identified TCL compound.
- E = Indicates that the compound was detected beyond the calibration range and was subsequently analyzed at a dilution.
- D = Identifies all compounds identified in an analysis at a secondary dilution factor.
- I = Interference.
- NQ = Result qualitatively confirmed but not able to quantify.
- A = Indicates that a TIC is a suspected aldol-condensation product.
- N = Indicates presumptive evidence of a compound. This flag is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It is applied to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the N code is not used.
- X = This flag is used for a TIC compound which is quantified relative to a response factor generated from a daily calibration standard (rather than quantified relative to the closest internal standard).
- Y = Additional qualifiers used as required are explained in the case narrative.

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IJ ,...

# **GLOSSARY OF BNA DATA**

### **ABBREVIATIONS**

BS = Indicates blank spike in which reagent grade water is spiked with the CLP matrix spike solutions and carried through all the steps in the method. Spike recoveries are reported.

BSD = Indicates blank spike duplicate.

MS = Indicates matrix spike.

MSD = Indicates matrix spike duplicate.

DL = Suffix added to sample number to indicate that results are from a diluted analysis.

NA = Not Applicable.

**DF** = Dilution Factor.

NR = Not Required.

SP, Z = Indicates Spiked Compound.

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Recra LabNet Philadelphia Sample Discrepancy Report (SDR) SDR#: 99 からつつ
Initiator: CTaylor RFW Batch: 99081851 Parameter: BNA Date: 91299 Samples: all Matrix: Water/cont Client: 1741-Hanford Method: SW846/MCAWW/CLP/ Prep Batch: 971E1035-ms
1. Reason for SDR a. COC Discrepancy Tech Profile Error Client Request Sampler Error on C-O-C Transcription Error Wrong Test Code Other b. General Discrepancy Missing Sample/Extract Container Broken Wrong Sample Pulled Label ID's lilegible Hold Time Exceeded Insufficient Sample Preservation Wrong Received Past Hold Improper Bottle Type Not Amenable to Analysis  Note: Verified by [Log-In] or [Prep Group] (circle)signature/date: C. QC Problem (Include all relevant specific results; attach data if necessary)  That Contained SX CRQ L of his (2 Extraplery)  Mthalate (65 Mb/L)
2. Known or Probable Causes(s)  Lab contamination.
3. Discussion and Proposed Action  Re-log  Entire Batch Following Samples: Re-leach Re-extract Re-digest Revise EDD Change Test Code to Place On/Take Off Hold (circle)  Other Description:  Sumple & QC were contained  Sumple & QC were contained  The Description:  The Description:  Sumple & QC were
4. Project Manager Instructionssignature/date:  Concur with Proposed Action Disagree with Proposed Action; See Instruction Include in Case Narrative Client Contacted: Date/Person Add Cancel
5. Final Actionsignature/date: CT 9 M/PS Other Explanation:  Verified re-[log][leach][extract][digest][analysis] (circle)  Included in Case Narrative  Hard Copy COC Revised  Electronic COC Revised  EDD Corrections Completed
When Final Action has been recorded, forward original to QA Specialist for distribution and filing.
Route Distribution of Completed SDR Route Distribution of Completed SDR  X Initiator Metals: Doughty  X Lab Manager: M. Taylor Inorganic: Perrone  X Project Mgr: Stone/Carey/Schrenkel/Johnson  X Section Mgr: Wesson/Daniels GC/LC: Schnell  X Section Mgr: Wesson/Daniels MS: LeMin/Taylor  X QA (file): Racioppi Log-in: Toder  Data Management: Feldman Admin: Soos  Sample Prep: Schnell/Doughty/Kauffman Other:

# Recra LabNet - Lionville Laboratory

Semivolatiles by GC/MS, HSL List

Report Date: 09/28/99 16:12 💆 Client: TNU-HANFORD B99-085 Work Order: 10985001001 Page: 1a RFW Batch Number: 9908L851

	Cust ID:	B0W67	9	B0W67	9	B0W679	)	SBLKBT		SBLKBT BS		
Sample Information	RFW#: Matrix:	001 WATER	L	001 MS WATER	3	001 MSD WATER	)	99LE1035-M WATER	B1	99LE1035-M WATER	181.	
1111011114111111	D.F.:	1.0	0.0	1.(	10	1.0	0	1.0	0	1.0	0	
	Units:	UG/I		UG/I		UG/L	-	UG/L		UG/L		
	5.11c5.	00,1	_	33, -	_			, _		•		
	Nitrobenzene-d5	78	%	79	%	81	%	82	%	65	%	
Surrogate	2-Fluorobiphenyl	72	%	80	%	73	8	72	જ	55	%	
Recovery	Terphenyl-d14	97	%	96	웅	87	8	95	૪	78	%	
	Phenol-d5	59	8	59	%	50	%	55	8	50	%	
	2-Fluorophenol	50	ક	52	%	52	%	51	왐	46	8	
	2,4,6-Tribromophenol	69	96	74	ે	77	%	69	8	58	8	
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Phenol		12	U	63	8	55	%	10	U	49	%	
bis(2-Chloro	ethyl)ether	12	U	24	U	24	U	10	U	10	U	
2-Chloropheno	01	12		71	%	64	%		U	51	%	
1,3-Dichlorobenzene		12		24	U	24	Ü	10	U	10	Ū	
1,4-Dichlorob	benzene	12		74	8	69	%	- •	U	59	ૄ	
	benzene	12		24	U	24	U	10	Ü	10	U	
2-Methylpheno			U	24	U	24	U	10	U	10	U	
2,2'-oxybis(	1-Chloropropane)	12		24	U	24	U	10	U	10	U	
	ol	12	U	24	Ū	24	Ü	10	U	10	U	
	-n-propylamine		U	81	જ	98	૪	10	U	81	ક	
	hane	12	_	24	U	24	U	10	U	10	U	
		12		24	U	24	U	10	Ü	10	U	
		12	U	24	U	24	U	10	U	10	U	
	1	12	U	24	U	24	Ŭ	10	U	10	_	
	phenol	12	U	24	U	24	U	10	U	10	U	
	ethoxy) methane	12	U	24	U	24	Ū	10	U	10 10	U	
	phenol	12		24	Ü	24	U	10	IJ	61	٠ ا	
	orobenzene	12	U	70	왕	64	8	10	U	10	5 U	
Naphthalene_		12		24	U. U	24	U	10 10	IJ	10	U	
	ine	12		24	-	24	U	10	U	10	Ū	
	tadiene		U	24	U •.	24	∪ %		U	5 <b>4</b>	\ %	
	ethylphenol	12	U	60	8	60	-	10	U	10	Ü	
2-Methylnaphthalene		12	U	24	U	24	U	10	U	10	U	
	clopentadiene	12		24	U	24	U	10	U		Ü	
2,4,6-Trichlo		12	U	24	U	24	U	10	_	10 25	U	
2,4,5-Trichlo		30	U	60	U	60	U	25	U	25	U	
*= Outside of EPA CLP QC limits.												

Fluoranthene\_\_\_\_\_

Indeno(1,2,3-cd)pyrene\_\_\_\_

Pyrene\_\_\_\_\_

Butylbenzylphthalate\_\_\_\_\_

24 U

24 U

24 U

24 U

왐

86

24 U

24 U

24 U

24 U

%

77

10 U

1.0 U

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74

3,3'-Dichlorobenzidine\_\_\_\_ 12 U 24 U 24 U 10 U 10 U Benzo(a)anthracene\_\_\_\_\_ 12 U 24 U 24 U 10 U 10 U Chrysene 12 U 24 U 24 U 10 U 10 U bis(2-Ethylhexyl)phthalate\_\_\_\_ 51 B 59 B 10 JB 65 24 В Di-n-octyl phthalate\_\_\_\_\_ 12 U 24 ·U 24 U 10 U 10 IJ Benzo(b)fluoranthene\_\_\_\_ 12 U 24 U 24 U 10 U 10 U Benzo(k)fluoranthene\_\_\_\_\_ 12 U 24 U 24 II 10 [] 10 IJ Benzo(a)pyrene\_\_\_\_ 12 U 24 U 24 U 10 U 10 U

Dibenz(a,h)anthracene\_\_\_\_\_ Benzo(g,h,i)perylene\_\_\_\_\_ 12 U 24 U 24 U 10 U (1) - Cannot be separated from Diphenylamine. \*= Outside of EPA CLP QC limits.

12 U

12 U

12 U

12 U

12 U

CLIENT SAMPLE NO.

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

| |B0W679 |

Lab Name: <u>Recra.LabNet</u>

Work Order: 10985001001

Client: TNU-HANFORD B99-085

Matrix: (soil/water) WATER Lab Sample ID: 9908L851-001

Sample wt/vol: 820 (g/mL)  $\underline{\text{ML}}$  Lab File ID:  $\underline{\text{A091312}}$ 

Level: (low/med) LOW Date Received: 08/24/99

% Moisture: \_\_\_\_ decanted: (Y/N)\_\_ Date Extracted: 08/26/99

Concentrated Extract Volume: 1000(uL) Date Analyzed: 09/13/99

Injection Volume: 2.0(uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: 7.0

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=======================================	= =====================================	=====	==========	=====
1.	UNKNOWN	7.16	2	J
2.	UNKNOWN	7.96	2	J
3.	UNKNOWN	8.23	4	J
4.	UNKNOWN	8.49	4	J
5.	UNKNOWN	8.59	5	J
6.	UNKNOWN	23.72	5	J
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CLIENT SAMPLE NO.

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET
TENTATIVELY IDENTIFIED COMPOUNDS

1			
SBLKBT	•		

Lab Name: <u>Recra.LabNet</u>

Work Order: 10985001001

Client: TNU-HANFORD B99-085

Matrix: (soil/water) WATER Lab Sample ID: 99LE1035-MB1

Sample wt/vol:  $\underline{1000}$  (g/mL)  $\underline{ML}$  Lab File ID:  $\underline{A091307}$ 

Level: (low/med) LOW Date Received: 08/26/99

% Moisture: \_\_\_\_ decanted: (Y/N)\_ Date Extracted: 08/26/99

Concentrated Extract Volume: 1000(uL) Date Analyzed: 09/13/99

Injection Volume: 2.0(uL) Dilution Factor: 1.00

GPC Cleanup: (Y/N) N pH: 7.0

CONCENTRATION UNITS:

Number TICs found: 8 (ug/L or ug/Kg) UG/L

}				
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
=======================================	=======================================	======		=====
1.	UNKNOWN	7,27	5	J
2.	UNKNOWN	7.96	10	J
3.	UNKNOWN	8.02	4	J
4.	UNKNOWN	8.23	6	J
5. ·	UNKNOWN	8.26	3	J
6.	UNKNOWN	8.50	6	J
7.	UNKNOWN	8.58	10	J
8.	UNKNOWN	8.66	2	J
	1			

# Recra LabNet - Lionville Laboratory BNA ANALYTICAL DATA PACKAGE FOR TNU-HANFORD B99-085

DATE RECEIVED: 08/24/99 RFW LOT # :9908L851

CLIENT ID	RFW #	MTX	PREP #	COLLECTION	EXTR/PREP	ANALYSIS
B0W679	001	W	99LE1035	08/19/99	08/26/99	09/13/99
B0W679	001 MS	M	99LE1035	08/19/99	08/26/99	09/13/99
B0W679	001 MSD	W	99LE1035	08/19/99	08/26/99	09/13/99
LAB QC:						
<u>·</u>						
SBLKBT	MB1	W	99LE1035	N/A	08/26/99	09/13/99
SBLKBT	MB1 BS	W	99LE1035	N/A	08/26/99	09/13/99

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<b>OL -</b> Drum Liquids		-						+		<del>                                     </del>	<del> </del>		<del> </del>											-	
L- EP/TCLP Leachate		<del>                                     </del>					+			<del>                                     </del>	<u> </u>		<del>                                     </del>	_					,					1	
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8/27/99 added Be to metals per client 3. improperly yreserved bottle Airbill * * 2) Unbroken on Oute Package (Y) or N  metal = As, ba, Cd, Cr, Cw, Pb, Ti, Se, 2) Ambient or Chilled 3) Present on Sample  3) Present on Sample												or N													
COMPOSITE											U, 21			سب	<b>, LJ</b>	, 10		<u>-</u> ,	(3)	Receive	ed in G	ood	3) Pres		Sample or N
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